

# WEST Search History

Hide Items

Restore

Clear

Cancel

DATE: Tuesday, March 02, 2004

Hide?	Set Name	Query	Hit Count
		<i>DB=PGPB,USPT,USOC,EPAB,JPAB,DWPI; PLUR=YES; OP=ADJ</i>	
<input type="checkbox"/>	L14	5213811.pn. or 4663318.pn.	4
<input type="checkbox"/>	L13	L12 and tablet	17
<input type="checkbox"/>	L12	L11 and l10	30
<input type="checkbox"/>	L11	particle	1358117
<input type="checkbox"/>	L10	L9 and l8	57
<input type="checkbox"/>	L9	controlled release	35915
<input type="checkbox"/>	L8	l1 or l2 or l4 or l5	375
<input type="checkbox"/>	L7	lycoremine	0
<input type="checkbox"/>	L6	lycoremin	0
<input type="checkbox"/>	L5	jilkon	1
<input type="checkbox"/>	L4	galantamin	1
<input type="checkbox"/>	L3	galantamina	0
<input type="checkbox"/>	L2	galanthamine	293
<input type="checkbox"/>	L1	galantamine	99

END OF SEARCH HISTORY

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(FILE 'HOME' ENTERED AT 17:31:14 ON 02 MAR 2004)

FILE 'REGISTRY' ENTERED AT 17:31:30 ON 02 MAR 2004

L1 1 S GALANTAMINE/CN

FILE 'USPATFULL, CAPLUS, MEDLINE, BIOSIS' ENTERED AT 17:34:51 ON 02 MAR 2004

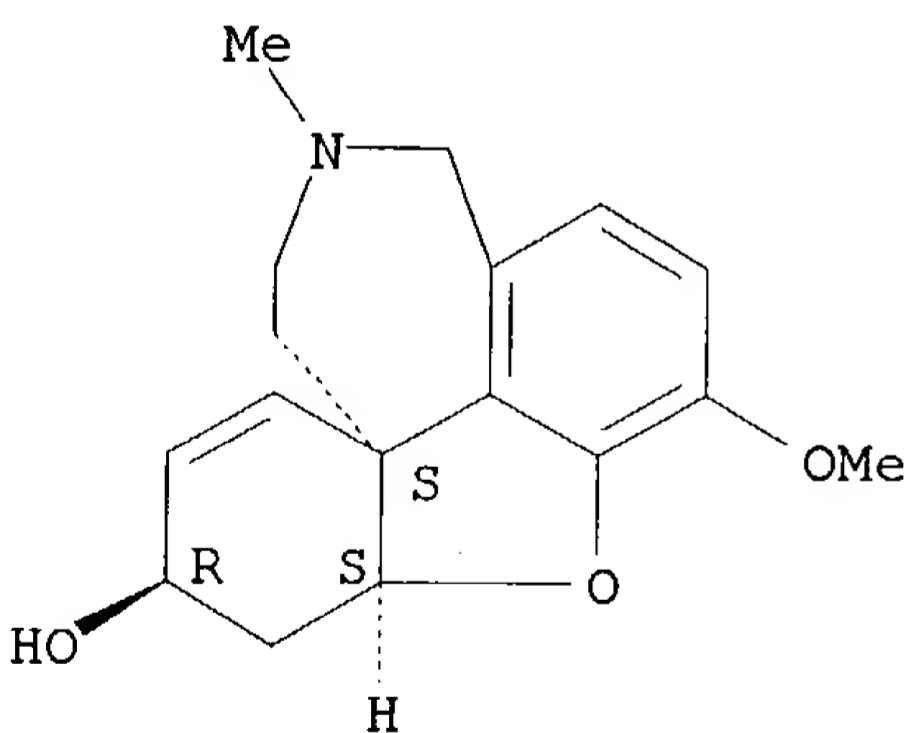
L2 1742 S L1  
L3 50281 S CONTROLLED RELEASE  
L4 1768319 S PARTICLES OR PARTICLE  
L5 1833135 S COAT?  
L6 16 S L2 AND L3  
L7 11 S L2 AND L4  
L8 29 S L2 AND L5  
L9 8 S L8 AND L4  
L10 6 S L8 AND L3  
L11 2 S L10 AND L9 AND L7 AND L6  
L12 16 DUP REM L6 (0 DUPLICATES REMOVED)  
L13 11 DUP REM L7 (0 DUPLICATES REMOVED)  
L14 6 S L8 AND L3

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L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN  
RN 357-70-0 REGISTRY  
CN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, (4aS,6R,8aS)- (9CI) (CA INDEX NAME)  
OTHER CA INDEX NAMES:  
CN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl- (7CI)  
CN Galanthamine (6CI, 8CI)  
OTHER NAMES:  
CN (-)-Galanthamine  
CN 6H-Benzofuro[3a,3,2-ef][2]benzazepin-6-ol, 4a,5,9,10,11,12-hexahydro-3-methoxy-11-methyl-, [4aS-(4a $\alpha$ ,6 $\beta$ ,8aR\*)]-  
CN BRN 0093736  
CN Galantamin  
CN Galantamina  
CN **Galantamine**  
CN Jilkon  
CN Lycoremin  
CN Lycoremine  
CN NSC 100058  
CN [4aS-(4a $\alpha$ ,6 $\beta$ ,8aR\*)]-4a,5,9,10,11,12-Hexahydro-3-methoxy-11-methyl-6H-benzofuro[3a,3,2-ef][2]benzazepin-6-ol  
FS STEREOSEARCH  
DR 736-79-8, 1551-02-6  
MF C17 H21 N O3  
CI COM  
LC STN Files: ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BEILSTEIN\*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CBNB, CEN, CHEMCATS, CIN, DDFU, DRUGU, EMBASE, HODOC\*, IMSDRUGNEWS, IMSPATENTS, IMSRESEARCH, IPA, MEDLINE, MRCK\*, NAPRALERT, PHAR, PROMT, RTECS\*, SPECINFO, SYNTHLINE, TOXCENTER, USAN, USPAT2, USPATFULL  
(\*File contains numerically searchable property data)  
Other Sources: WHO

Absolute stereochemistry. Rotation (-).



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

663 REFERENCES IN FILE CA (1907 TO DATE)  
22 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA  
665 REFERENCES IN FILE CAPLUS (1907 TO DATE)  
27 REFERENCES IN FILE CAOLD (PRIOR TO 1967)

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